

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE101618\
 Data File : BE097912.D
 Acq On : 16 Oct 2018 15:42
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE101618

Quant Time: Oct 16 17:37:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 15:36:44 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	101	0.00
2	1,4-Dioxane	0.400	0.384	4.0	105	0.00
3	n-Nitrosodimethylamine	0.400	0.380	5.0	108	0.00
4 S	2-Fluorophenol	0.400	0.385	3.8	103	0.00
5 S	Phenol-d6	0.400	0.390	2.5	101	0.00
6	bis(2-Chloroethyl)ether	0.400	0.382	4.5	103	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	105	0.00
8 S	Nitrobenzene-d5	0.400	0.402	-0.5	107	0.00
9	Naphthalene	0.400	0.394	1.5	106	-0.01
10	Hexachlorobutadiene	0.400	0.388	3.0	105	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	106	0.00
12	2-Methylnaphthalene	0.400	0.398	0.5	108	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	108	0.00
14 S	2,4,6-Tribromophenol	0.400	0.359	10.3	110	0.00
15 S	2-Fluorobiphenyl	0.400	0.379	5.3	108	0.00
16	Acenaphthylene	0.400	0.380	5.0	108	0.00
17	Acenaphthene	0.400	0.387	3.3	110	0.00
18	Fluorene	0.400	0.391	2.3	111	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	118	0.00
20	4-Bromophenyl-phenylether	0.400	0.376	6.0	113	0.00
21	Hexachlorobenzene	0.400	0.383	4.3	115	0.00
22	Pentachlorophenol	0.400	0.364	9.0	116	0.00
23	Phenanthrene	0.400	0.384	4.0	117	0.00
24	Anthracene	0.400	0.377	5.8	116	0.00
25 SURR	Fluoranthene-d10	0.400	0.380	5.0	120	0.00
26	Fluoranthene	0.400	0.378	5.5	119	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	106	0.00
28	Pvrene	0.400	0.418	-4.5	117	0.00
29 S	Terphenyl-d14	0.400	0.396	1.0	115	0.00
30	Benzo(a)anthracene	0.400	0.389	2.8	109	0.00
31	Chrysene	0.400	0.375	6.3	108	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.324	19.0	93	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.342	14.5	98	0.00
34 I	Perylene-d12	0.400	0.400	0.0	96	0.00
35	Benzo(b)fluoranthene	0.400	0.371	7.3	96	0.00
36	Benzo(k)fluoranthene	0.400	0.376	6.0	100	0.00
37 C	Benzo(a)pyrene	0.400	0.383	4.3	98	0.00
38	Dibenzo(a,h)anthracene	0.400	0.384	4.0	99	0.00
39	Benzo(a,h,i)perylene	0.400	0.379	5.3	100	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0				